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REILLY TAR & CHEMICAL CORPORATION

TELEPHONE 317-638-7531
CABLE RETAR INDIANAPOLIS
TELEX 27-404



1510 MARKET SQUARE CENTER
151 NORTH DELAWARE STREET
INDIANAPOLIS, INDIANA 46204

July 23, 1981

Illinois EPA
Division of Land/Noise Pollution Control
2200 Churchill Road
Springfield, Illinois 62706

Gentlemen:

Attached is our application for Special Waste. In place of the hauler registration number we are supplying their ICC number which is 15493-MC-C. Also attached is a copy of analysis on the waste material.

Very truly yours,

REILLY TAR & CHEMICAL CORPORATION

W. A. Justin
Director Environmental Control

WAJ/bk

Attach.

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STATE OF ILLINOIS

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ENVIRODYNE
engineers
2151 Lackland Road
St. Louis, Missouri 63141
314 424 6960

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REPORT OF ANALYSIS

Environment • Energy • Transportation • Food Processing

SUBMITTED BY: Mr. Gary Bush
c/o Enviro Chem
865 South State Road (421)
Zionxville, Indiana 46077

DATE: July 9, 1981
PROJECT NO. 1702-126
P.O.

DATE RECEIVED: June 18, 1981

SAMPLE ANALYZED: Coal Tar Residue

METHODS USED: Toxic Extraction Procedure, Federal Register, May 19, 1980;
EPA's Sludge Protocol for VOAs; EPA's Region V Methods; and EPA's
Hazardous Waste Manual, Method 810

RESULTS

<u>Leachate</u>		<u>Sample</u>	
Arsenic	0.021	% Volatile at 100°C	10%
Barium	0.12	% Volatile at 600°C	70%
Chromium	0.053	Flash Point	Did not flash, went up to 80°C
Cadmium	0.016	Cyanide	250
Lead	<0.05	Sulfide	7
Mercury	0.0010	pH	7.4 pH units
Silver	<0.02	TCDD	ND (<0.1)
Selenium	<0.02	Benzene	30.3
Endrin	<0.001	Toluene	18.2
Lindane	<0.001	phenol	* 125
Methoxychlor	<0.001	Carbazole**	250
Toxaphene	<0.001	Phenanthrene**	25,200
2,4-D	<0.001	Naphthalene**	14,300
2,4,5-TP	<0.001	Ethylbenzene	1.93
		Methylene chloride	5.41

Values are ppm except as otherwise noted.

GC/MS Library Search results for volatile identities are attached.

*Phenol results are not yet complete due to a backordered part.

**EPA Hazardous Waste Manual defines cresote as a combination of these compounds.

/csq

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ENVIRODYNE ENGINEERS

BY: *Paul Myers*

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1702-126
GC/MS LIBRARY SEARCH
VOLATILES

Peak No.	Retention Time	Similarity Index	Identification	Concentration (ppm)	Concentration in Blank** (ppm)
1	2.3	0.9778 0.7350	dimethyl ether* d -dimethylamine	1.65	5.5
2	7.0	-	acetone	1.92	1.8
3	12.4	0.9695 0.9691	1,2-diethoxyethane diethyl ether*	1.07	
4	13.7	-	methyl ethyl ketone	3.0	4.2
5	14.9	-	pentane	1.82	0.93
6	16.2	(No hits - not enough detail in spectrum)		0.585	
7	21.3	-	hexane	0.768	0.27
8	29.5	0.9807 0.9806 0.9804 0.9784	1-methylphenylacetylene* indene* benzene, 1-ethynyl-4-methyl-* benzene, 1,3-propadienyl-	276	

* = best match

**Concentrations in blank calculated to compare with sample. All peaks except No. 8 are probably background and not really present in the sample.

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